

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
 Data File : BM042869.D
 Acq On : 18 Nov 2023 07:02
 Operator : MA/JU
 Sample : PB156976BL
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK976

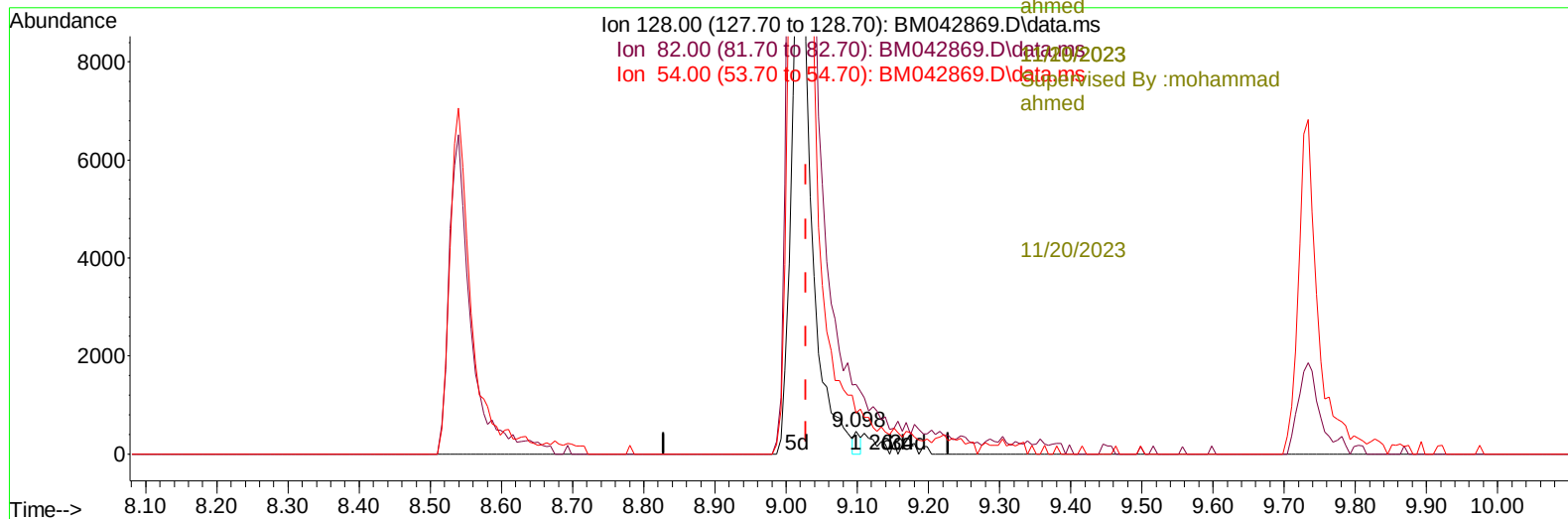
Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/19/2023
 Supervised By :mohammad ahmed 11/20/2023

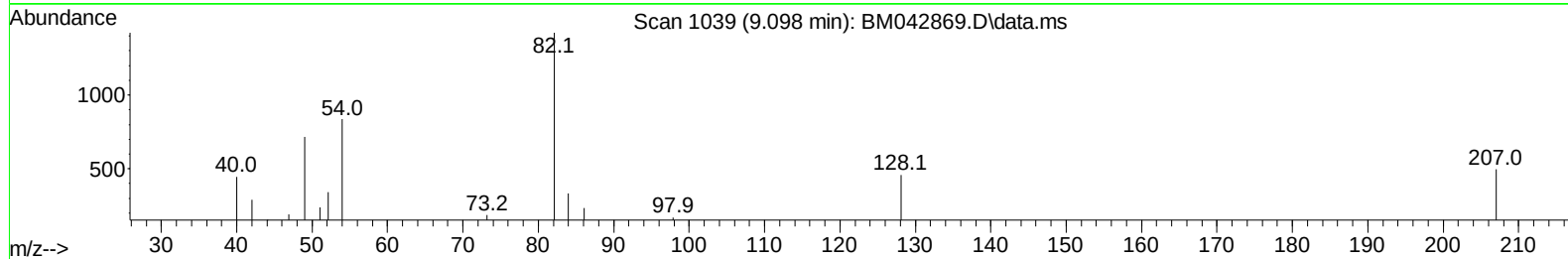
11/19/2023
 Supervised By :mohammad
 ahmed
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Quant Time: Nov 18 21:06:55 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 15 22:28:07 2023
 Response via : Initial Calibration

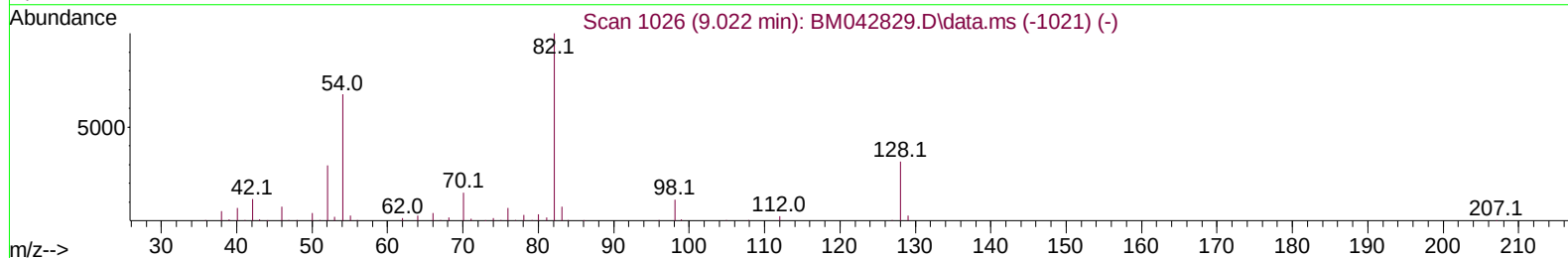
Ion 128.00 (127.70 to 128.70): BM042869.D\data.ms
 Ion 82.00 (81.70 to 82.70): BM042869.D\data.ms
 Ion 54.00 (53.70 to 54.70): BM042869.D\data.ms



Scan 1039 (9.098 min): BM042869.D\data.ms



Scan 1026 (9.022 min): BM042829.D\data.ms (-1021) (-)



TIC: BM042869.D\data.ms

(21) Nitrobenzene-d5 (S)

9.098min (+ 0.070) 0.32 ng/uL

response	281	
Ion	Exp%	Act%
128.00	100.00	100.00
82.00	289.30	309.80
54.00	185.60	181.92
0.00	0.00	0.00

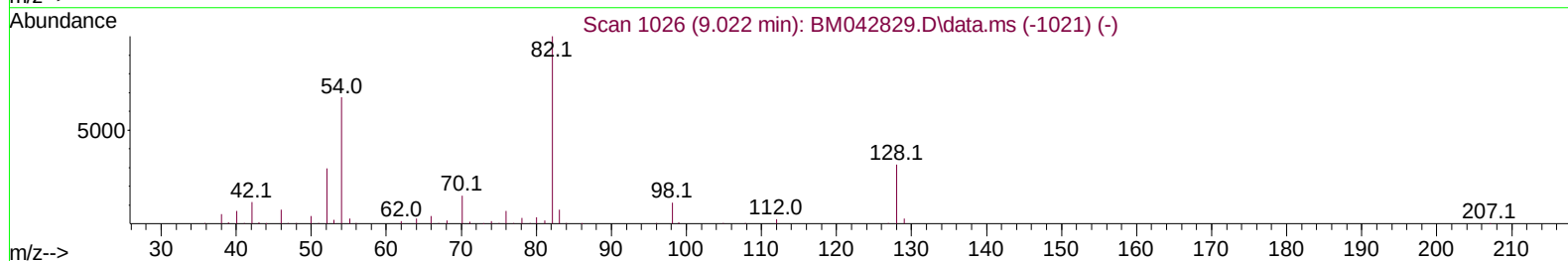
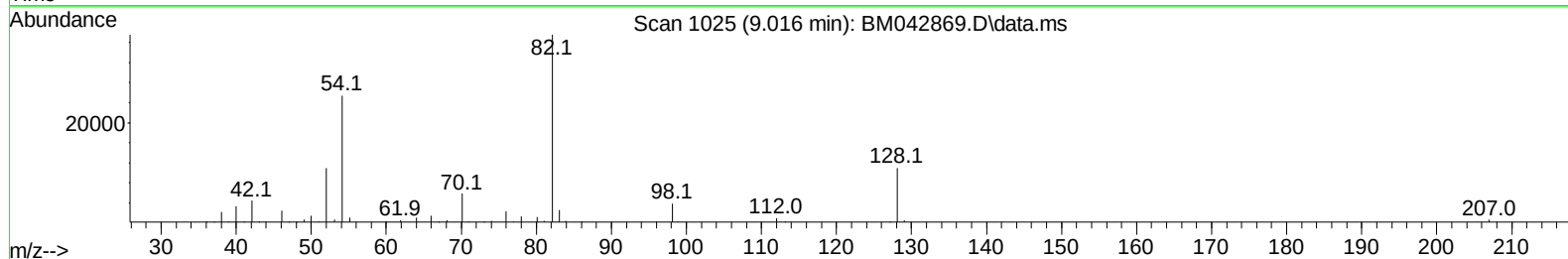
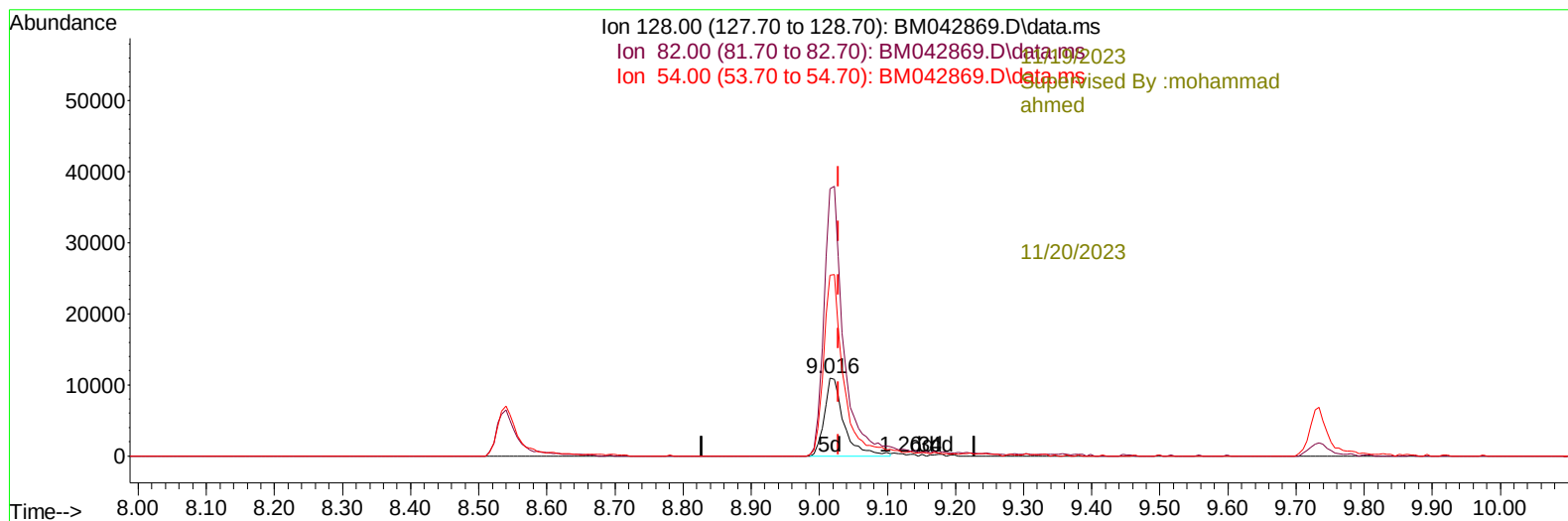
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TIC: BM042869.D\data.ms

(21) Nitrobenzene-d5 (S)

9.016min (-0.012) 24.55 ng/ul m

response 21717

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	289.30	342.23
54.00	185.60	231.87#
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

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Quant Time: Nov 18 21:25:05 2023
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)
							-11/19/2023
							Supervised By :mohammad
							ahmed
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.816	152		25543	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.628	136		100897	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.474	164		62726	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.227	188		134332	20.000	ng/ul	-0.01
79) Chrysene-d12	21.415	240		131202	20.000	ng/ul	0.00
88) Perylene-d12	23.768	264		166066	20.000	ng/ul	-0.02
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.175	96		5072	6.231	ng/uL	0.00
4) Pyridine-d5	3.610	84		57915	28.602	ng/ul	0.00
7) Phenol-d5	6.987	99		63284	25.456	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.163	67		53438	30.591	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.340	132		50704	26.433	ng/ul	-0.01
15) 4-Methylphenol-d8	8.540	113		46829	23.924	ng/ul	-0.01
21) Nitrobenzene-d5	9.016	128		21717m	24.550	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.734	143		23443	22.654	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.257	165		39918	20.745	ng/ul	0.00
31) 4-Chloroaniline-d4	10.810	131		52518	21.496	ng/ul	0.00
46) Dimethylphthalate-d6	13.898	166		159204	26.817	ng/ul	0.00
49) Acenaphthylene-d8	14.174	160		171192	26.951	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.751	143		11342m	16.910	ng/ul	0.01
60) Fluorene-d10	15.469	176		135630	27.590	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.621	200		15755	15.484	ng/ul	0.00
73) Anthracene-d10	17.327	188		202971	26.961	ng/ul	-0.01
81) Pyrene-d10	19.621	212		277159	32.259	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.615	264		285688	28.760	ng/ul	-0.01

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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